

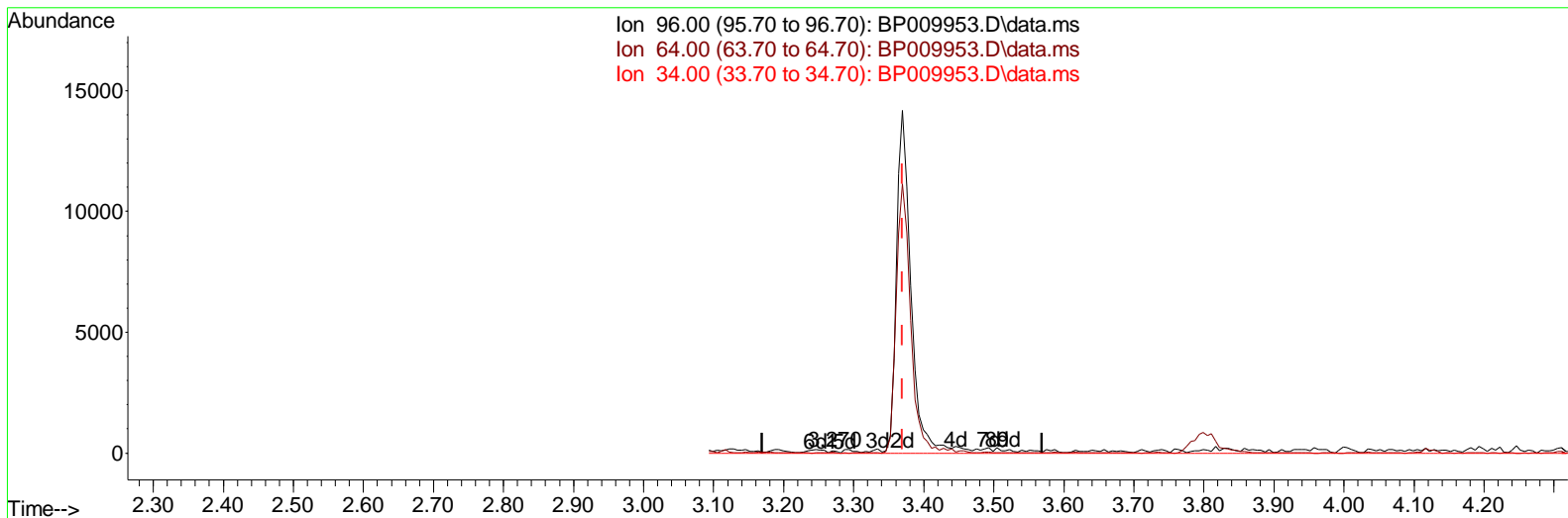
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 Data File : BP009953.D
 Acq On : 19 Apr 2022 19:44
 Operator : CG/JU
 Sample : PB144172BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLCS172

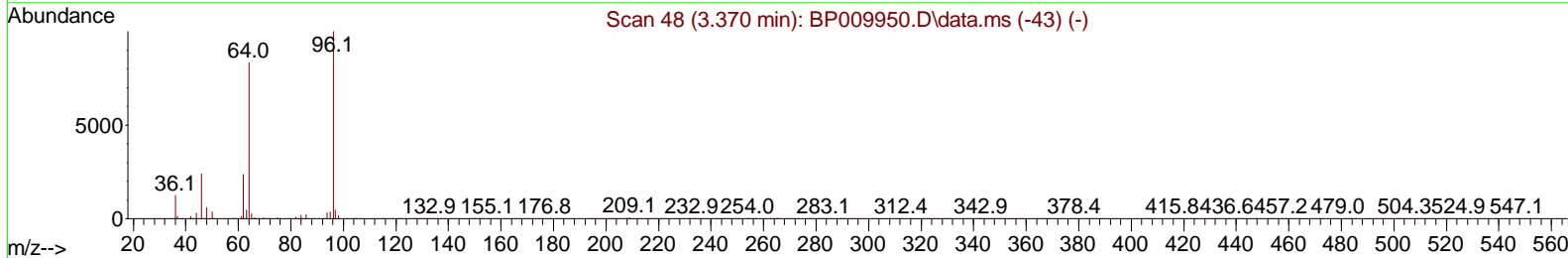
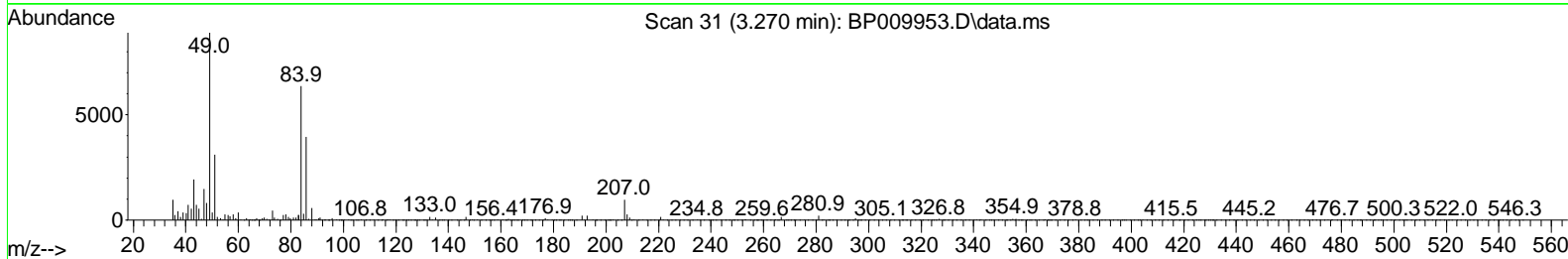
Manual Integrations APPROVED

Reviewed By : Jagrut Upadhyay 04/20/2022
 Supervised By : Yogesh Patel 04/25/2022

Quant Time: Apr 20 01:33:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 20 01:32:14 2022
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Ion 96.00 (95.70 to 96.70): BP009953.D\data.ms
 Ion 64.00 (63.70 to 64.70): BP009953.D\data.ms
 Ion 34.00 (33.70 to 34.70): BP009953.D\data.ms



TIC: BP009953.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.270min (-0.100) 0.01 ng/uL

| response | 47 | |
|----------|--------|--------|
| Ion | Exp% | Act% |
| 96.00 | 100.00 | 100.00 |
| 64.00 | 55.40 | 45.98 |
| 34.00 | 0.00 | 0.00 |
| 0.00 | 0.00 | 0.00 |

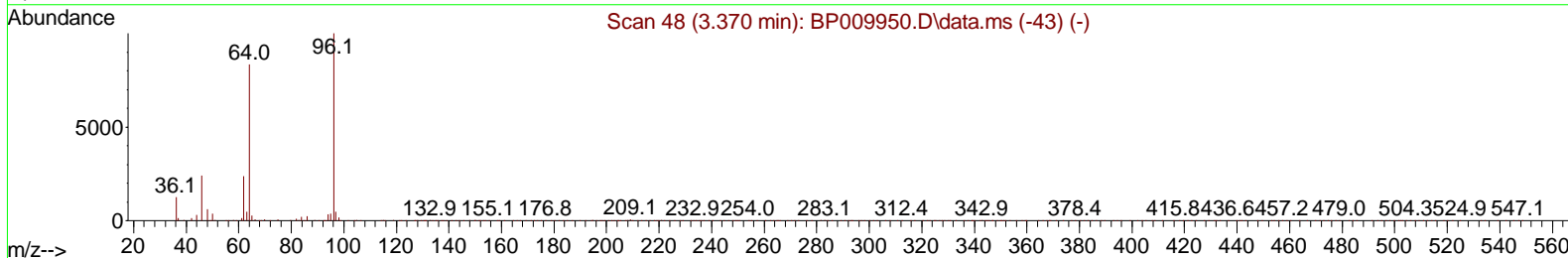
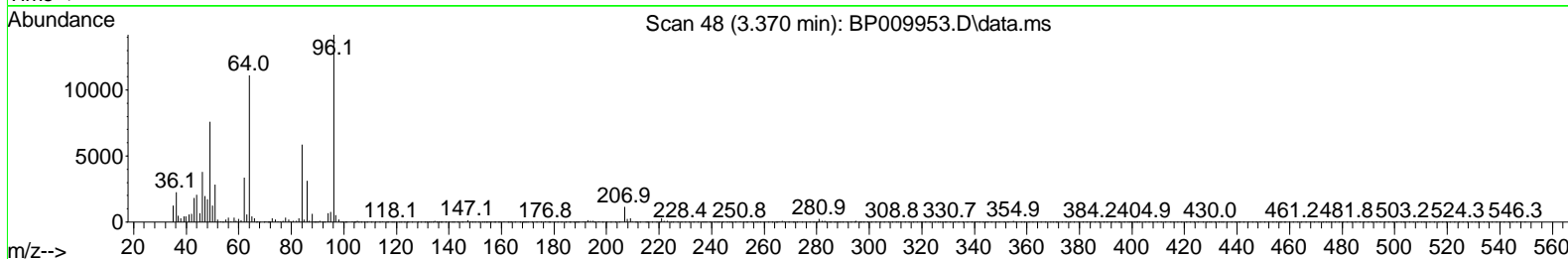
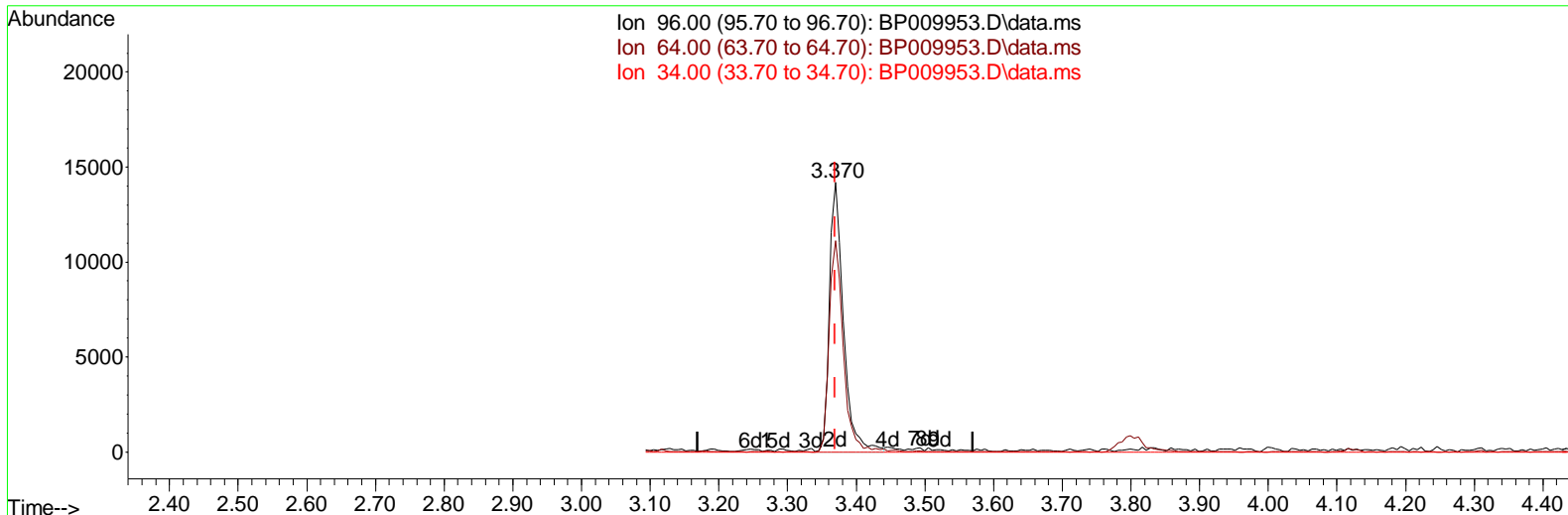
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 Sample : PB144172BS
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 ALS Vial : 5 Sample Multiplier: 1

Instrument :
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 Client Sample Id :
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TIC: BP009953.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.370min (-0.000) 5.98 ng/uL m

| response | 20118 | |
|----------|--------|--------|
| Ion | Exp% | Act% |
| 96.00 | 100.00 | 100.00 |
| 64.00 | 55.40 | 78.47# |
| 34.00 | 0.00 | 0.00 |
| 0.00 | 0.00 | 0.00 |

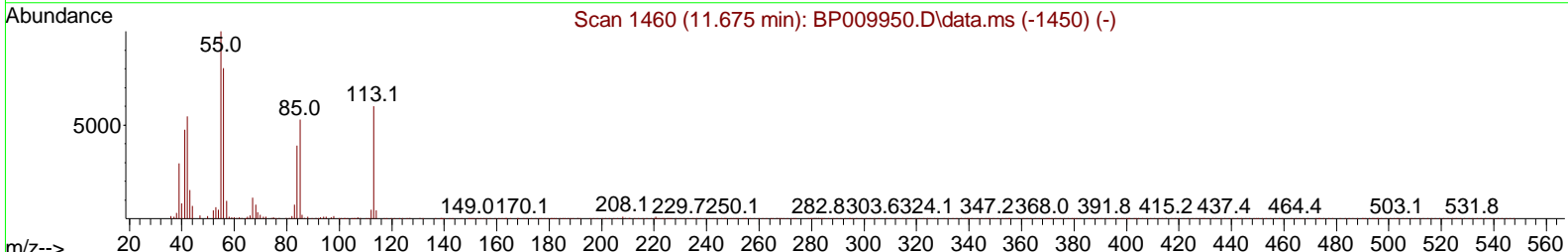
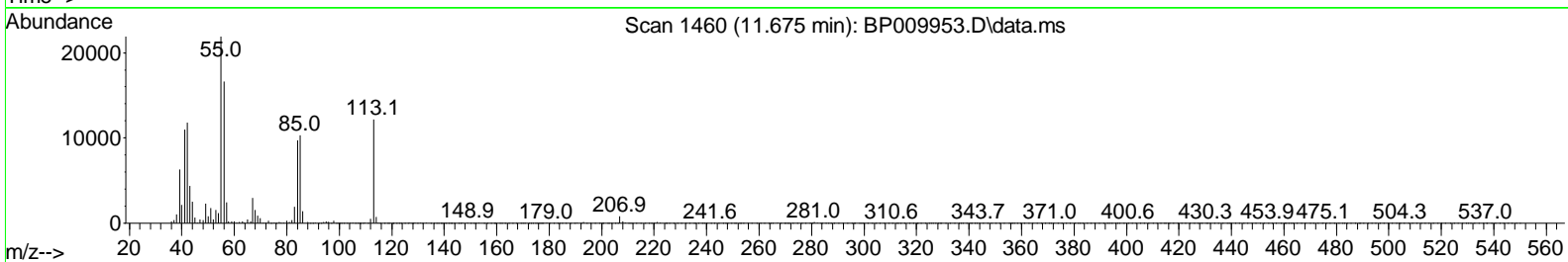
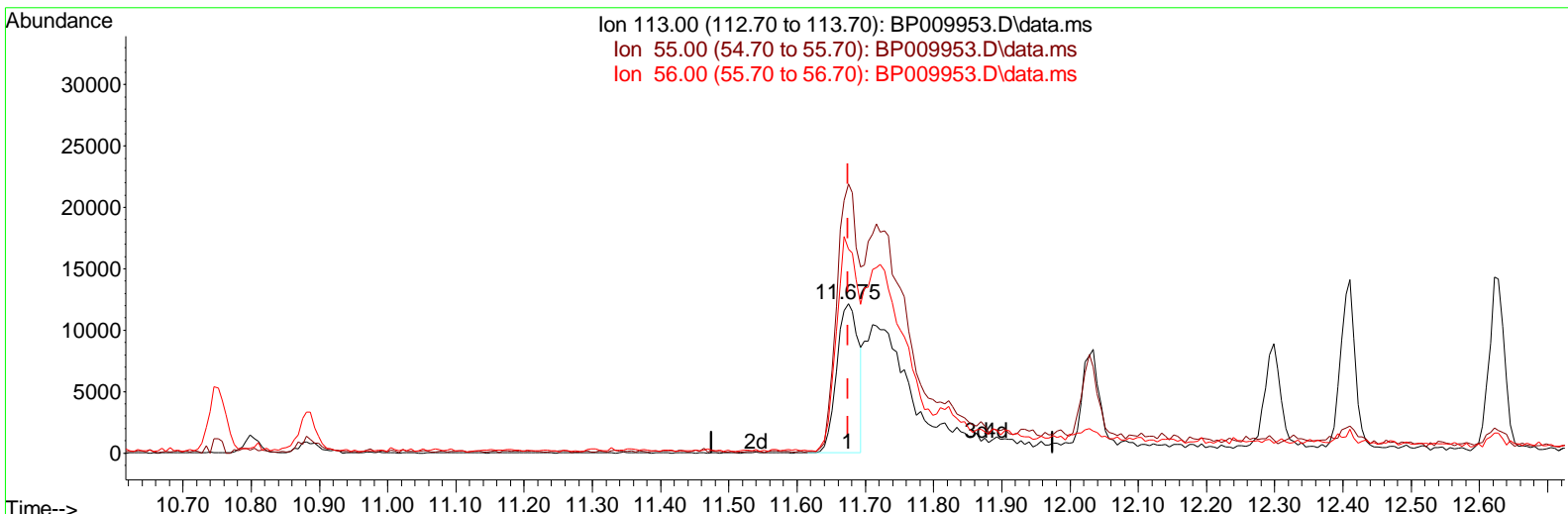
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
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TIC: BP009953.D\data.ms

(34) Caprolactam

11.675min (-0.000) 7.56 ng/ul

response 26849

| Ion | Exp% | Act% |
|--------|--------|---------|
| 113.00 | 100.00 | 100.00 |
| 55.00 | 72.70 | 179.83# |
| 56.00 | 85.80 | 136.47# |
| 0.00 | 0.00 | 0.00 |

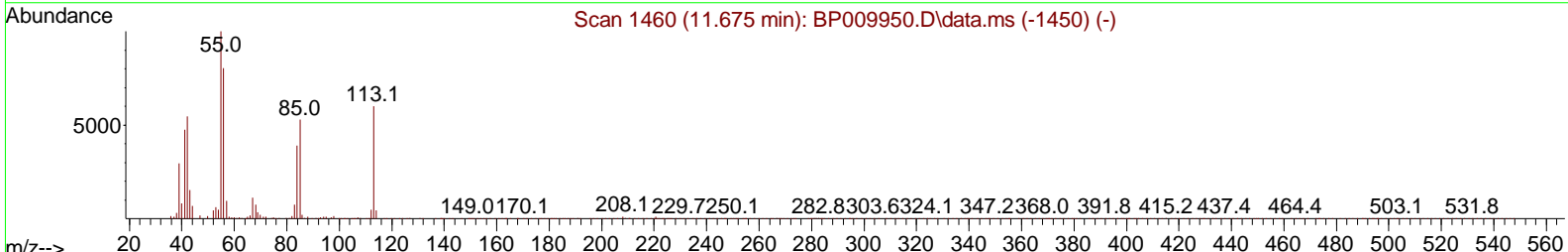
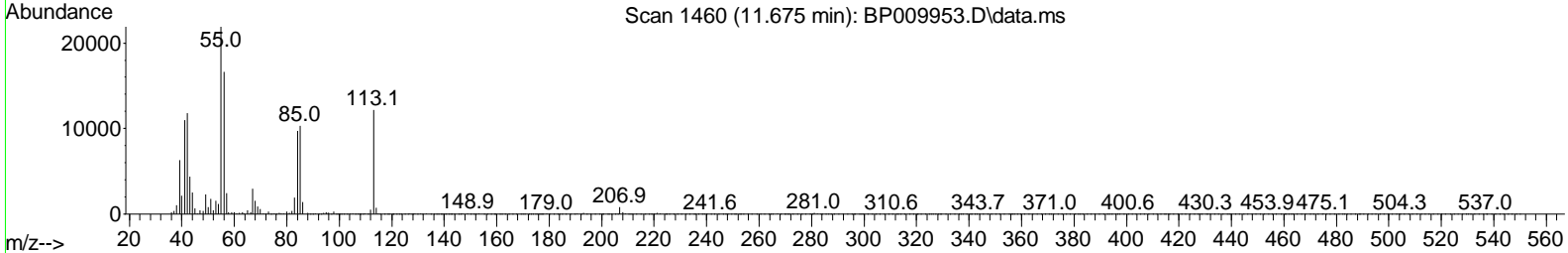
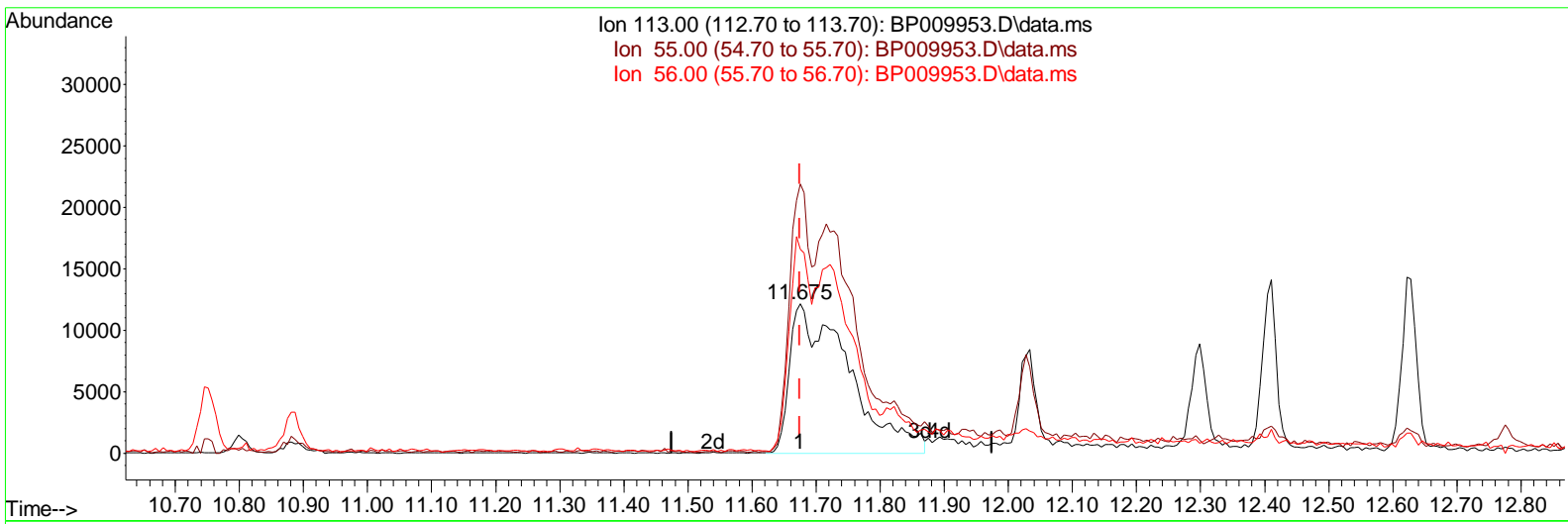
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TIC: BP009953.D\data.ms

(34) Caprolactam

11.675min (-0.000) 21.88 ng/ul m

response 77673

| Ion | Exp% | Act% |
|--------|--------|---------|
| 113.00 | 100.00 | 100.00 |
| 55.00 | 72.70 | 179.83# |
| 56.00 | 85.80 | 136.47# |
| 0.00 | 0.00 | 0.00 |

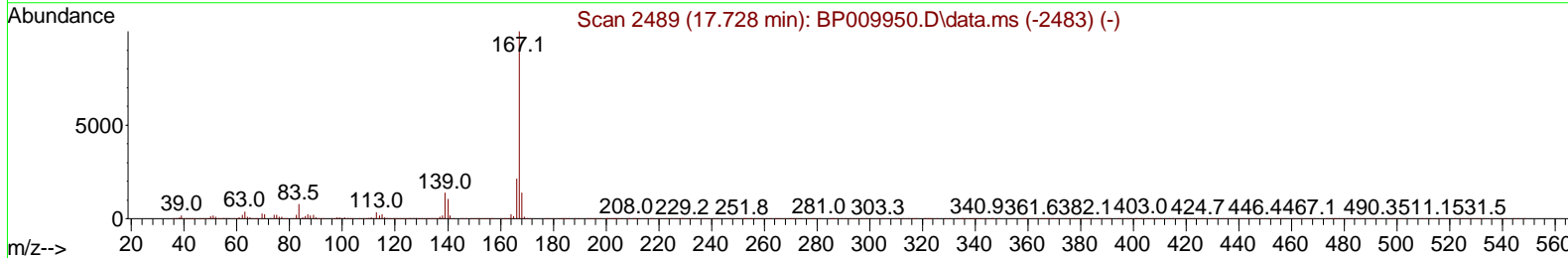
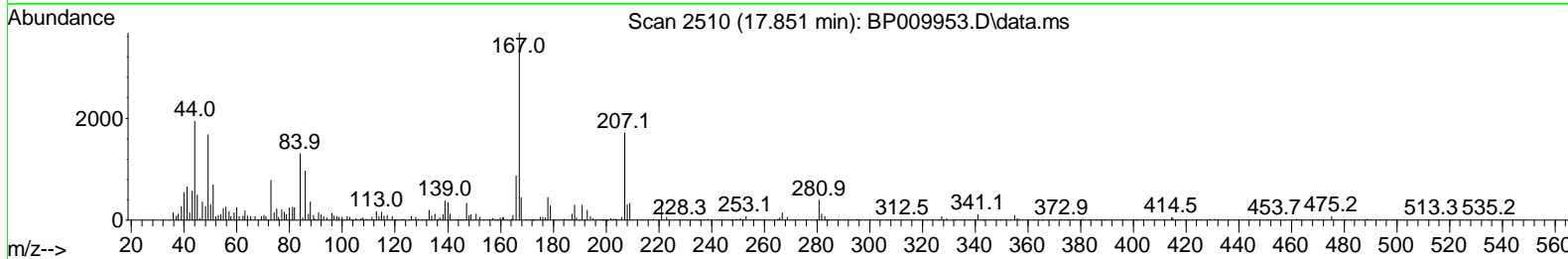
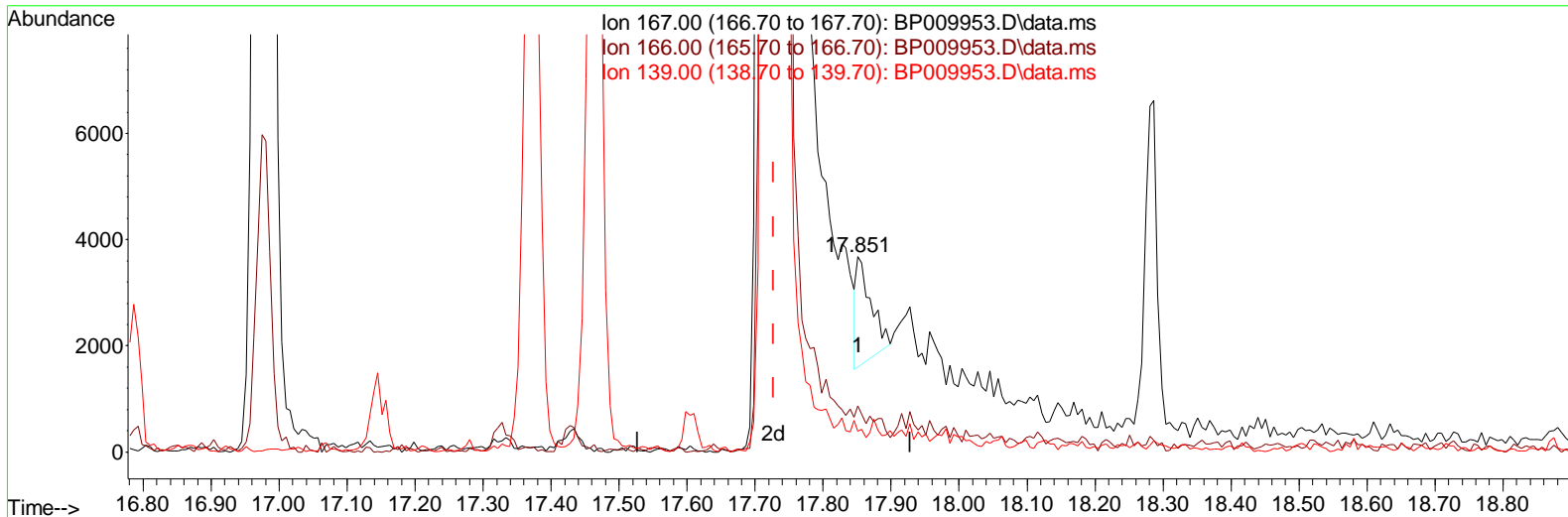
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TIC: BP009953.D\data.ms

(77) Carbazole

17.851min (+ 0.123) 0.06 ng/ul

response 3075

| Ion | Exp% | Act% |
|--------|--------|--------|
| 167.00 | 100.00 | 100.00 |
| 166.00 | 19.80 | 23.61 |
| 139.00 | 10.40 | 10.52 |
| 0.00 | 0.00 | 0.00 |

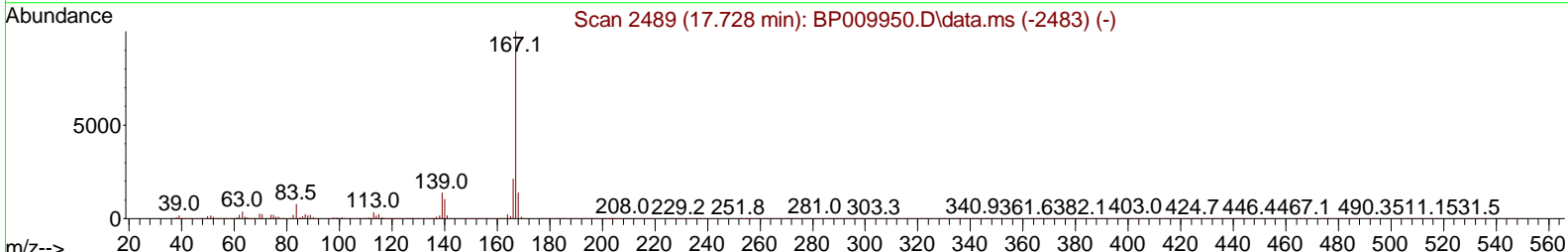
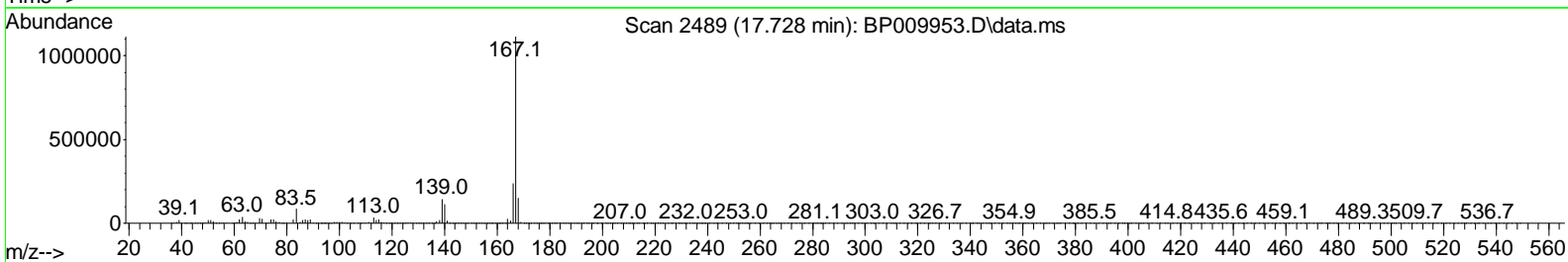
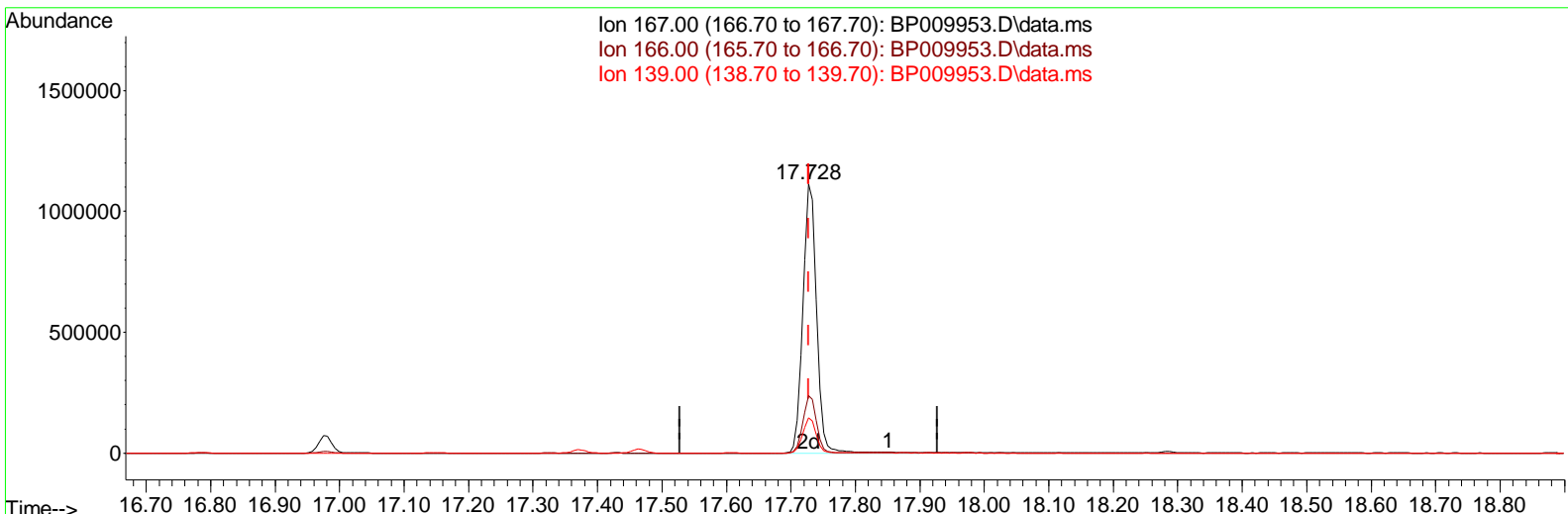
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Instrument :
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TIC: BP009953.D\data.ms

(77) Carbazole

17.728min (-0.000) 32.32 ng/ul m

response 1648381

| Ion | Exp% | Act% |
|--------|--------|--------|
| 167.00 | 100.00 | 100.00 |
| 166.00 | 19.80 | 21.35 |
| 139.00 | 10.40 | 13.01# |
| 0.00 | 0.00 | 0.00 |

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 Misc :
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Instrument :
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ClientSampleId :
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 Response via : Initial Calibration

| Compound | R. T. | QI on | Response | Conc | Units | Dev(Min) |
|------------------------------------|--------|-------|----------|--------|---------|----------|
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 7.946 | 152 | 150101 | 20.000 | ng/ul | 0.00 |
| 20) Naphthalene-d8 | 10.752 | 136 | 669239 | 20.000 | ng/ul | 0.00 |
| 38) Acenaphthene-d10 | 14.575 | 164 | 472105 | 20.000 | ng/ul | 0.00 |
| 64) Phenanthrene-d10 | 17.328 | 188 | 1042692 | 20.000 | ng/ul | # 0.00 |
| 79) Chrysene-d12 | 21.410 | 240 | 1038282 | 20.000 | ng/ul | # 0.00 |
| 88) Perylene-d12 | 23.868 | 264 | 995084 | 20.000 | ng/ul | 0.00 |
| System Monitoring Compounds | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.370 | 96 | 20118m | 5.982 | ng/uL | 0.00 |
| 4) Pyridine-d5 | 3.781 | 84 | 264326 | 28.505 | ng/ul | 0.00 |
| 7) Phenol-d5 | 7.105 | 99 | 397072 | 30.132 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)eth... | 7.269 | 67 | 253160 | 32.269 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4 | 7.475 | 132 | 320750 | 32.268 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8 | 8.652 | 113 | 344711 | 31.335 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5 | 9.105 | 128 | 167028 | 34.609 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4 | 9.828 | 143 | 188587 | 35.642 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3 | 10.369 | 165 | 349277 | 31.546 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4 | 10.881 | 131 | 427245 | 27.093 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6 | 13.987 | 166 | 1188620 | 32.052 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8 | 14.269 | 160 | 1342652 | 30.560 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4 | 14.769 | 143 | 207826 | 30.901 | ng/ul | 0.00 |
| 60) Fluorene-d10 | 15.569 | 176 | 1018448 | 32.646 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.687 | 200 | 212733 | 34.440 | ng/ul | 0.00 |
| 73) Anthracene-d10 | 17.428 | 188 | 1600321 | 32.056 | ng/ul | 0.00 |
| 81) Pyrene-d10 | 19.657 | 212 | 1940351 | 34.218 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12 | 23.710 | 264 | 1708669 | 32.973 | ng/ul | 0.00 |
| Target Compounds | | | | | | |
| 2) 1,4-Dioxane | 3.405 | 88 | 39428 | 11.500 | ng/uL# | 18 |
| 5) Pyridine | 3.799 | 79 | 279771 | 28.917 | ng/ul # | 39 |
| 6) Benzaldehyde | 7.075 | 77 | 252511 | 31.244 | ng/ul | 95 |
| 8) Phenol | 7.134 | 94 | 424355 | 32.171 | ng/ul # | 95 |
| 10) Bis(2-Chloroethyl)ether | 7.363 | 93 | 330350 | 32.151 | ng/ul # | 77 |
| 12) 2-Chlorophenol | 7.505 | 128 | 327367 | 32.666 | ng/ul | 94 |
| 13) 2-Methylphenol | 8.387 | 108 | 325481 | 31.643 | ng/ul | 97 |
| 14) 2,2'-oxybis(1-Chloropr... | 8.481 | 45 | 484479 | 31.899 | ng/ul # | 86 |
| 16) Acetophenone | 8.769 | 105 | 533517 | 30.782 | ng/ul # | 81 |
| 17) N-Nitrosodipropylamine | 8.758 | 70 | 284470 | 31.658 | ng/ul # | 72 |
| 18) 4-Methylphenol | 8.716 | 108 | 353886 | 31.770 | ng/ul | 92 |
| 19) Hexachloroethane | 9.034 | 117 | 134713 | 31.838 | ng/ul | 85 |
| 22) Nitrobenzene | 9.146 | 77 | 412238 | 33.279 | ng/ul # | 84 |
| 23) Isophorone | 9.675 | 82 | 805270 | 31.547 | ng/ul | 96 |
| 25) 2-Nitrophenol | 9.857 | 139 | 198226 | 35.126 | ng/ul # | 86 |
| 26) 2,4-Dimethylphenol | 9.922 | 107 | 421469 | 31.472 | ng/ul # | 84 |
| 27) Bis(2-Chloroethoxy)meth... | 10.157 | 93 | 474608 | 31.669 | ng/ul # | 97 |
| 29) 2,4-Dichlorophenol | 10.399 | 162 | 348655 | 32.481 | ng/ul # | 82 |
| 30) Naphthalene | 10.799 | 128 | 1101827 | 32.012 | ng/ul | 96 |
| 32) 4-Chloroaniline | 10.904 | 127 | 430687 | 27.599 | ng/ul | 98 |
| 33) Hexachlorobutadiene | 11.099 | 225 | 245492 | 31.755 | ng/ul | 94 |
| 34) Caprolactam | 11.675 | 113 | 77673m | 21.881 | ng/ul | |
| 35) 4-Chloro-3-methylphenol | 12.028 | 107 | 414878 | 31.898 | ng/ul # | 71 |

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 Response via : Ini tial Cal i brati on

| Compound | R. T. | QI on | Response | Conc | Units | Dev(Mi n) |
|-------------------------------------|--------|-------|----------|--------|---------|-----------|
| 36) 2-Methyl naphthal ene | 12.410 | 142 | 794120 | 31.581 | ng/ul | 92 |
| 37) 1-Methyl naphthal ene | 12.628 | 142 | 806147 | 31.751 | ng/ul # | 95 |
| 39) 1, 2, 4, 5-Tetrachl oroben. . . | 12.775 | 216 | 466113 | 32.270 | ng/ul # | 96 |
| 40) Hexachl orocycl opentadi ene | 12.763 | 237 | 294327 | 29.364 | ng/ul | 96 |
| 41) 2, 4, 6-Tri chl orophenol | 13.010 | 196 | 274948 | 29.807 | ng/ul # | 81 |
| 42) 2, 4, 5-Tri chl orophenol | 13.081 | 196 | 286615 | 28.752 | ng/ul # | 87 |
| 43) 1, 1' -Bi phenyl | 13.410 | 154 | 1122911 | 32.599 | ng/ul | 93 |
| 44) 2-Chl oronaphthal ene | 13.457 | 162 | 864873 | 32.663 | ng/ul | 93 |
| 45) 2-Ni troani li ne | 13.651 | 65 | 290335 | 37.416 | ng/ul # | 82 |
| 47) Di methyl phthal ate | 14.034 | 163 | 1158365 | 32.454 | ng/ul # | 93 |
| 48) 2, 6-Di ni trotol uene | 14.145 | 165 | 244279 | 37.208 | ng/ul | 96 |
| 50) Acenaphthyl ene | 14.298 | 152 | 1420872 | 32.771 | ng/ul | 95 |
| 51) 3-Ni troani li ne | 14.475 | 138 | 216361 | 31.700 | ng/ul # | 82 |
| 52) Acenaphthene | 14.640 | 153 | 914240 | 32.205 | ng/ul | 96 |
| 53) 2, 4-Di ni trophenol | 14.675 | 184 | 106427 | 26.818 | ng/ul # | 80 |
| 55) 4-Ni trophenol | 14.781 | 109 | 185794 | 30.010 | ng/ul # | 53 |
| 56) Di benzofuran | 14.975 | 168 | 1340706 | 32.089 | ng/ul | 99 |
| 57) 2, 4-Di ni trotol uene | 14.934 | 165 | 360587 | 37.404 | ng/ul # | 84 |
| 58) 2, 3, 4, 6-Tetrachl orophenol | 15.204 | 232 | 297967 | 32.783 | ng/ul # | 88 |
| 59) Di ethyl phthal ate | 15.398 | 149 | 1203314 | 32.576 | ng/ul | 93 |
| 61) Fl uorene | 15.628 | 166 | 1110925 | 32.281 | ng/ul | 91 |
| 62) 4-Chl orophenyl -phenyl e. . . | 15.622 | 204 | 590969 | 32.296 | ng/ul | 98 |
| 63) 4-Ni troani li ne | 15.639 | 138 | 240745 | 35.216 | ng/ul # | 80 |
| 66) 4, 6-Di ni tro-2-methyl ph. . . | 15.698 | 198 | 208354 | 33.935 | ng/ul # | 90 |
| 67) N-Ni trosodi phenyl ami ne | 15.834 | 169 | 986140 | 33.039 | ng/ul | 94 |
| 68) 4-Bromophenyl -phenyl ether | 16.516 | 248 | 365752 | 32.405 | ng/ul | 95 |
| 69) Hexachl orobenzene | 16.639 | 284 | 419140 | 32.295 | ng/ul # | 91 |
| 70) Atrazi ne | 16.786 | 200 | 400597 | 32.516 | ng/ul | 91 |
| 71) Pentachl orophenol | 16.981 | 266 | 245809 | 27.981 | ng/ul # | 84 |
| 72) Phenanthrene | 17.369 | 178 | 1825239 | 32.794 | ng/ul | 99 |
| 74) Anthracene | 17.463 | 178 | 1851360 | 32.809 | ng/ul | 98 |
| 75) 1, 2, 3, 4-Tetrachl oroben. . . | 13.381 | 216 | 496205 | 32.916 | ng/uL# | 86 |
| 76) Pentachl orobenzene | 14.898 | 250 | 477374 | 32.488 | ng/uL | 91 |
| 77) Carbazol e | 17.728 | 167 | 1648381m | 32.324 | ng/ul | |
| 78) Di -n-butyl phthal ate | 18.281 | 149 | 2107764 | 33.612 | ng/ul # | 93 |
| 80) Fl uoranthene | 19.333 | 202 | 2255047 | 34.549 | ng/ul # | 95 |
| 82) Pyrene | 19.686 | 202 | 2306236 | 34.850 | ng/ul # | 93 |
| 83) Butyl benzyl phthal ate | 20.557 | 149 | 968458 | 35.674 | ng/ul # | 79 |
| 84) 3, 3' -Di chl orobenzi di ne | 21.322 | 252 | 684402 | 30.933 | ng/ul # | 97 |
| 85) Benzo(a)anthracene | 21.398 | 228 | 2179904 | 33.118 | ng/ul | 95 |
| 86) Bi s(2-ethyl hexyl)phtha. . . | 21.322 | 149 | 1424422 | 34.883 | ng/ul # | 83 |
| 87) Chrysene | 21.451 | 228 | 2151093 | 33.872 | ng/ul | 99 |
| 89) Di -n-octyl phthal ate | 22.263 | 149 | 2417368 | 34.673 | ng/ul | 100 |
| 90) Benzo(b)fl uoranthene | 23.121 | 252 | 2244228 | 33.945 | ng/ul | 99 |
| 91) Benzo(k)fl uoranthene | 23.169 | 252 | 2049389 | 33.970 | ng/ul # | 98 |
| 93) Benzo(a)pyrene | 23.763 | 252 | 2068794 | 33.758 | ng/ul | 99 |
| 94) I ndeno(1, 2, 3-cd)pyrene | 26.439 | 276 | 2081450 | 32.480 | ng/ul # | 94 |
| 95) Di benzo(a, h)anthracene | 26.457 | 278 | 1765602 | 31.773 | ng/ul # | 96 |
| 96) Benzo(g, h, i)peryl ene | 27.221 | 276 | 1733898 | 32.814 | ng/ul # | 93 |

(#) = qual i fier out of range (m) = manual i ntegrati on (+) = signal s summed

Instrument :

BNA_P

ClientSampleId :

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